

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
 Data File : BP014401.D
 Acq On : 30 Mar 2023 14:45
 Operator : CG/JU
 Sample : 02061-01
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB961

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
 Title : SVOA CALIBRATION

Signal : TIC: BP014401.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.393	31	35	40	rVB	33203	43820	1.07%	0.109%
2	3.834	108	110	133	rBV	125732	256373	6.27%	0.639%
3	4.046	142	146	152	rVB	14129	24886	0.61%	0.062%
4	4.146	159	163	169	rVB	56223	80195	1.96%	0.200%
5	4.252	178	181	188	rVB7	8782	13864	0.34%	0.035%
6	4.858	281	284	289	rBV7	2249	4132	0.10%	0.010%
7	5.081	316	322	330	rBV2	11675	19249	0.47%	0.048%
8	5.270	350	354	364	rVB2	2615	6519	0.16%	0.016%
9	6.393	541	545	549	rBV3	6055	9064	0.22%	0.023%
10	7.181	673	679	693	rBV	114241	222418	5.44%	0.554%
11	7.340	700	706	718	rBV	527579	819789	20.04%	2.043%
12	7.546	733	741	751	rBV	471615	798811	19.52%	1.991%
13	8.016	814	821	831	rBV	628305	997851	24.39%	2.487%
14	8.575	912	916	919	rBV6	2719	4533	0.11%	0.011%
15	8.734	936	943	962	rBV	373175	675934	16.52%	1.685%
16	9.193	1014	1021	1037	rBV	643843	1098199	26.84%	2.737%
17	9.352	1043	1048	1053	rBV8	3760	6900	0.17%	0.017%
18	9.569	1079	1085	1092	rVB3	13934	26447	0.65%	0.066%
19	9.922	1137	1145	1160	rBV	509670	924185	22.59%	2.303%
20	10.463	1228	1237	1260	rBV	620372	1298169	31.73%	3.236%
21	10.610	1260	1262	1269	rVB8	5516	10749	0.26%	0.027%
22	10.840	1292	1301	1309	rBV	895977	1507736	36.85%	3.758%
23	10.987	1319	1326	1343	rBV	290568	649025	15.86%	1.618%
24	11.681	1441	1444	1445	rBV3	5043	4209	0.10%	0.010%
25	12.246	1533	1540	1544	rBV4	9904	17501	0.43%	0.044%
26	12.440	1568	1573	1581	rVB3	12017	23948	0.59%	0.060%
27	13.134	1688	1691	1696	rBV7	2192	4237	0.10%	0.011%
28	13.475	1744	1749	1754	rBV4	10001	15070	0.37%	0.038%
29	13.557	1757	1763	1765	rVV6	3021	4281	0.10%	0.011%
30	14.075	1843	1851	1866	rBV	1703336	2455259	60.01%	6.120%
31	14.181	1866	1869	1873	rVB5	4822	8307	0.20%	0.021%
32	14.363	1893	1900	1912	rBV	1808033	2638892	64.50%	6.577%
33	14.546	1925	1931	1934	rBV4	11363	13283	0.32%	0.033%
34	14.669	1945	1952	1961	rBV	1460762	2154468	52.66%	5.370%
35	14.928	1989	1996	2010	rBV3	36977	135912	3.32%	0.339%
36	15.493	2088	2092	2096	rVB3	15974	22605	0.55%	0.056%

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37	15.663	2115	2121	2129	rBV	2325183	3421614	83.63%	8.528%
38	15.793	2137	2143	2156	rBV	795861	1181046	28.87%	2.944%
39	15.910	2159	2163	2168	rVB4	16125	26504	0.65%	0.066%
40	16.034	2177	2184	2197	rBV	309951	454241	11.10%	1.132%
41	16.363	2235	2240	2248	rBV2	30851	54590	1.33%	0.136%
42	16.834	2316	2320	2321	rBV4	5118	6161	0.15%	0.015%
43	16.869	2324	2326	2330	rVB5	4765	6530	0.16%	0.016%
44	17.104	2363	2366	2372	rBV8	4716	6856	0.17%	0.017%
45	17.163	2372	2376	2380	rBV	20967	28577	0.70%	0.071%
46	17.434	2416	2422	2432	rBV	1736855	2532213	61.89%	6.311%
47	17.534	2432	2439	2452	rVV	2588150	3757248	91.83%	9.365%
48	17.704	2464	2468	2474	rVB	18992	27624	0.68%	0.069%
49	17.904	2499	2502	2506	rVB3	14217	15520	0.38%	0.039%
50	18.298	2564	2569	2576	rBV	80322	140874	3.44%	0.351%
51	18.369	2578	2581	2588	rVB	30453	50581	1.24%	0.126%
52	19.339	2742	2746	2752	rBV3	35577	49934	1.22%	0.124%
53	19.404	2754	2757	2766	rVB2	32608	65822	1.61%	0.164%
54	19.528	2775	2778	2784	rBV6	37344	61428	1.50%	0.153%
55	19.763	2812	2818	2825	rBV	2962666	4091351	100.00%	10.198%
56	21.198	3058	3062	3071	rBV3	110514	219398	5.36%	0.547%
57	21.522	3112	3117	3125	rVB	1422926	1998998	48.86%	4.982%
58	23.874	3510	3517	3527	rBV	1565796	3134568	76.61%	7.813%
59	24.039	3539	3545	3554	rVB2	843547	1792438	43.81%	4.468%

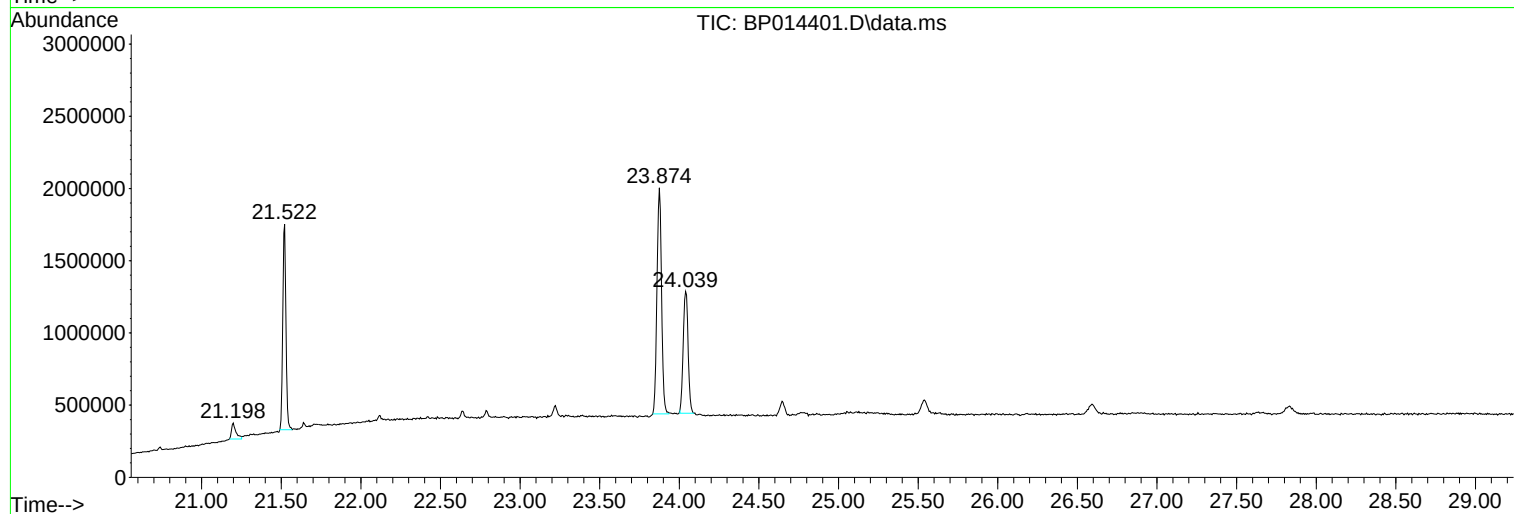
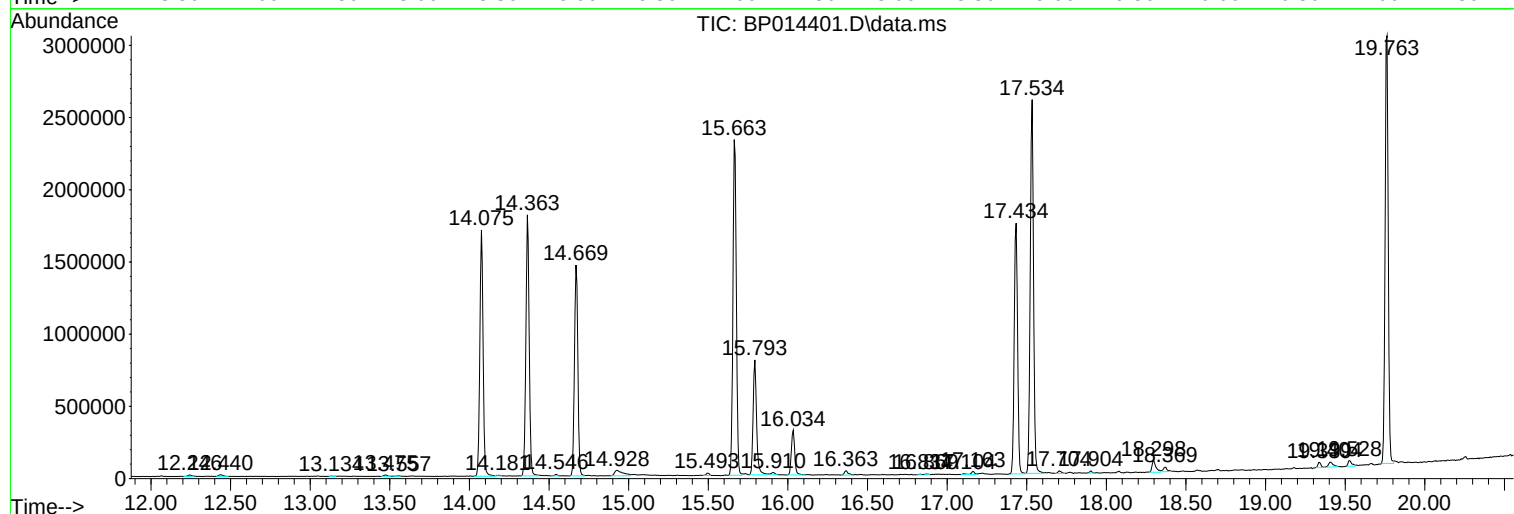
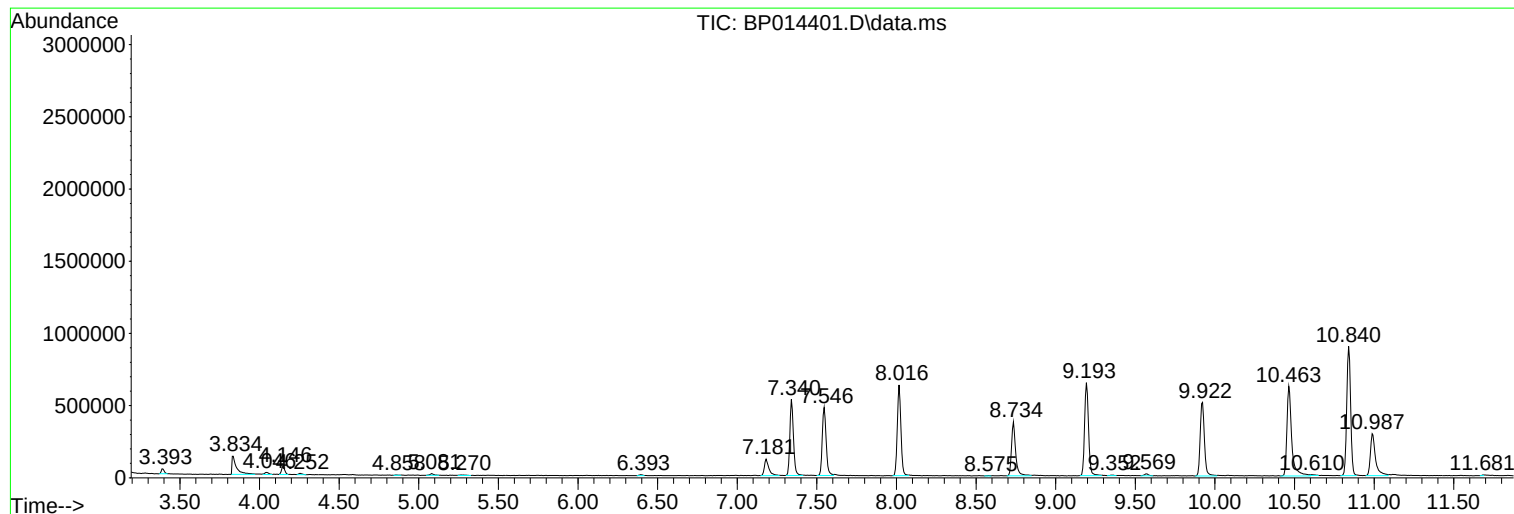
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



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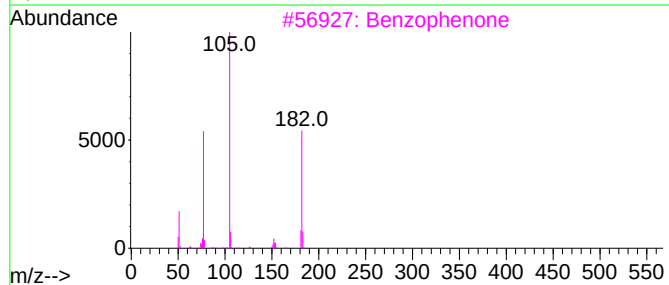
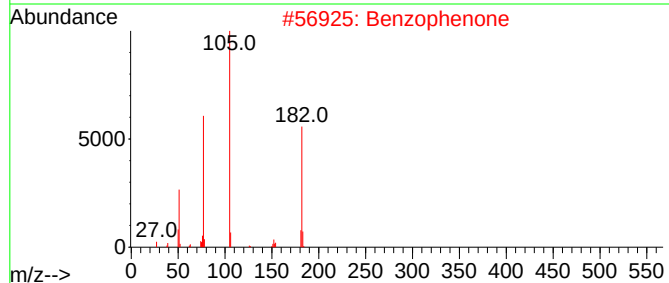
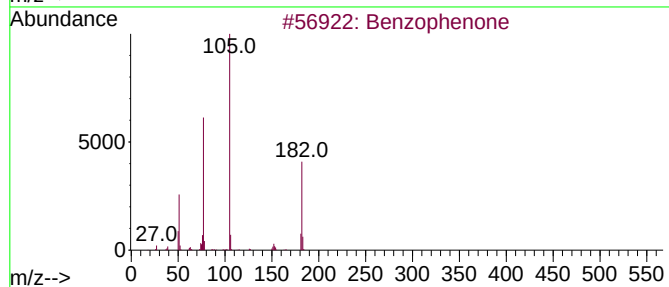
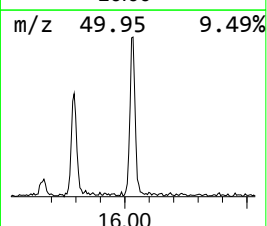
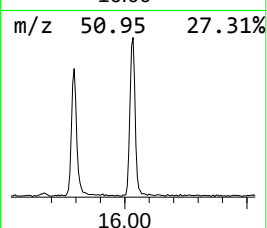
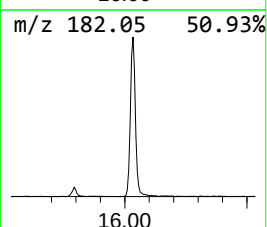
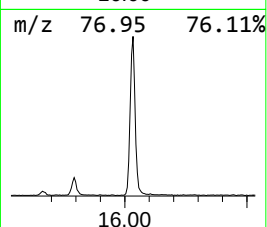
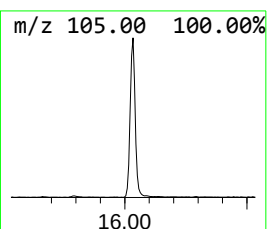
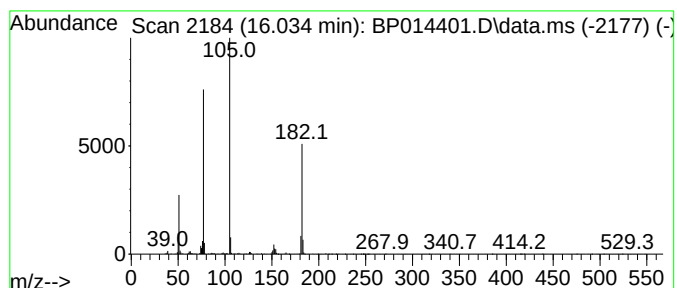
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Benzophenone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.034	4.22 ng/ul	454241	Acenaphthene-d10	14.669

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	95
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	87



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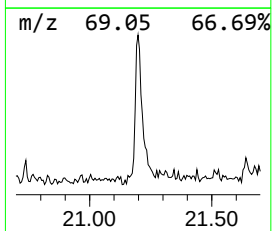
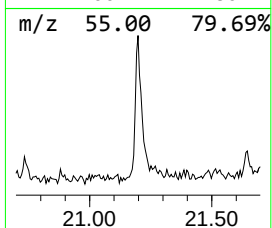
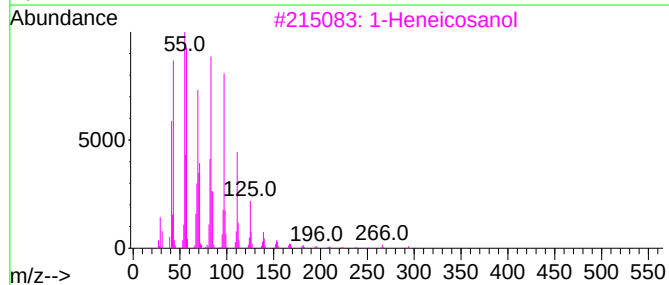
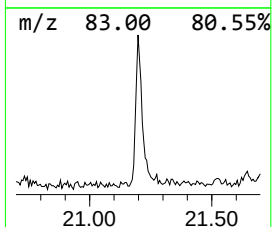
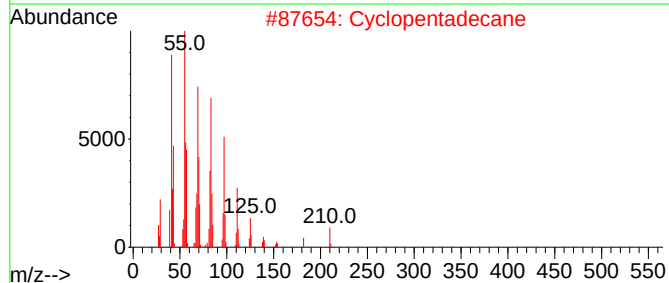
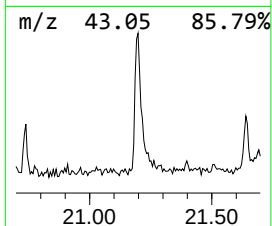
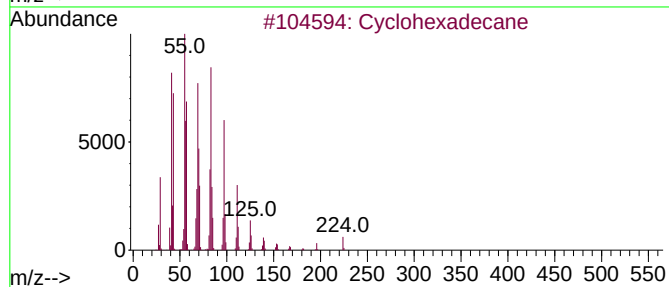
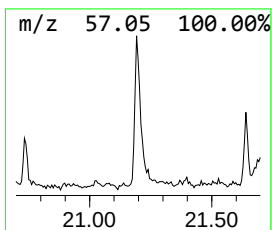
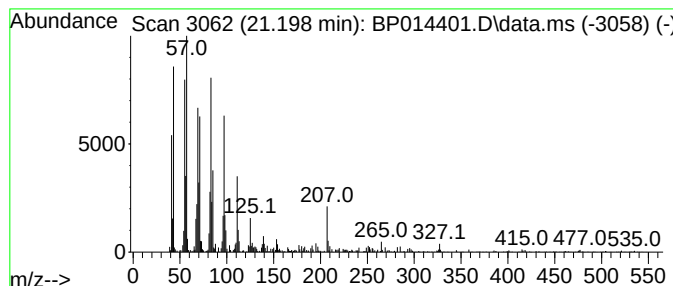
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 (DEL) Alkane: Cyclic21.198 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.198	2.20 ng/ul	219398	Chrysene-d12	21.522

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexadecane	224	C16H32	000295-65-8	94
2			Cyclopentadecane	210	C15H30	000295-48-7	93
3			1-Heneicosanol	312	C21H44O	015594-90-8	93
4			Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	91
5			n-Nonadecanol-1	284	C19H40O	001454-84-8	89



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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzophenone	16.034	4.2	ng/u1	454241	3	14.669	2154470	20.0
(DEL) Alkane: C...	21.198	2.2	ng/u1	219398	5	21.522	1999000	20.0